Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. (Currently Amended) A compound having the following formula (I):

$$\begin{array}{c|c} R_1 & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & \\ & & & \\ & & \\ & & & \\ & &$$

wherein

A is -(CHR3)- or -(C=O)-; ;

B is -(CHR₄)- or -(C=O)- $\frac{1}{2}$,

D is $-(CHR_5)$ - or -(C=O)- $\frac{1}{2}$,

E is $-(ZR_6)$ - or -(C=O)-;

G is (XR₇),

W is -Y(C=O)-, -(C=O)NH-, -(SO₂)- or nothing; 7

Y is oxygen, sulfur or -NH-; ;

X is nitrogen,

Z is nitrogen or CH; ; and

 $R_1,\,R_2,\,R_3,\,R_4,\,R_5,\,R_6,\,\text{and}\,\,R_7\,\,\text{are the same or different and independently selected from}\underline{:}$

an amino acid side chain moiety, an amino acid side chain derivative, a linker, and a solid support:

C1-12 alkyl, C6-12 aryl and C7-12 arylalkyl; and

 $\underline{aminoC_{2.5}alkyl}, \quad \underline{guanidinoC_{2.5}alkyl}, \quad \underline{C_{1.4}alkylguanidinoC_{2.5}alkyl}, \quad \underline{diC_{1.5}alkyl}$

 $_{\underline{4}}\underline{alkylguanidino} - \underline{C_{2.5}}\underline{alkyl}, \ \ \underline{amidino}\underline{C_{2.5}}\underline{alkyl}, \ \ \underline{C_{1.4}}\underline{alkylamidino}\underline{C_{2.5}}\underline{alkyl}, \ \ \underline{di}\underline{C_{1.4}}\underline{alkylamidino}\underline{C_{2.5}}\underline{alkyl},$

salkyl, C_{1-s}alkoxy, phenyl, substituted phenyl (where the substituents on the phenyl are independently selected from one or more of amino, amidino, guanidino, hydrazino, amidazonyl, C₁₋₄alkylamino, C₁₋₄dialkylamino, halogen, perfluoro C₁₋₄alkyl, C₁₋₄alkyl, C₁₋₃alkoxy, nitro, carboxy, cyano, sulfuryl or hydroxyl), benzyl, substituted benzyl (where the substituents on the benzyl are independently selected from one or more of amino, amidino, guanidino, hydrazino, amidazonyl, C1-4alkylamino, C1-4dialkylamino, halogen, perfluoro C1-4alkyl, C1-3alkoxy, nitro, carboxy, cyano, sulfuryl or hydroxyl), naphthyl, substituted naphthyl (where the substituents on the naphthyl are independently selected from one or more of amino, amidino, guanidino, hydrazino, amidazonyl, C₁₋₄alkylamino, C₁₋₄dialkylamino, halogen, perfluoro C₁₋₄alkyl, C₁-4alkyl, C1-3alkoxy, nitro, carboxy, cyano, sulfuryl or hydroxyl), bis-phenyl methyl, substituted bis-phenyl methyl (where the substituents on the bis-phenyl methyl are independently selected from one or more of amino, amidino, guanidino, hydrazino, amidazonyl, C₁₋₄alkylamino, C₁ 4dialkylamino, halogen, perfluoro C1-4alkyl, C1-4alkyl, C1-3alkoxy, nitro, carboxy, cyano, sulfuryl or hydroxyl), pyridyl, substituted pyridyl (where the substituents on the pyridyl are independently selected from one or more of amino amidino, guanidino, hydrazino, amidazonyl, C1-4alkylamino, C1-4dialkylamino, halogen, perfluoro C1-4alkyl, C1-4alkyl, C1-3alkoxy, nitro, carboxy, cyano, sulfuryl or hydroxyl), pyridylC₁₋₄alkyl, substituted pyridylC₁₋₄alkyl (where the pyridine substituents are independently selected from one or more of amino, amidino, guanidino, hydrazino, amidazonyl, C₁₋₄alkylamino, C₁₋₄dialkylamino, halogen, perfluoro C₁₋₄alkyl, C₁ 4alkyl, C₁₋₃alkoxy, nitro, carboxy, cyano, sulfuryl or hydroxyl), pyrimidylC₁₋₄alkyl, substituted pyrimidylC₁₋₄alkyl (where the pyrimidine substituents are independently selected from one or more of amino, amidino, guanidino, hydrazino, amidazonyl, C1-4alkylamino, C1-4dialkylamino, halogen, perfluoro C₁₋₄alkyl, C₁₋₄alkyl, C₁₋₃alkoxy, nitro, carboxy, cyano, sulfuryl or hydroxyl), triazin-2-yl-C₁₋₄alkyl, substituted triazin-2-yl-C₁₋₄alkyl (where the triazine substituents are independently selected from one or more of amino, amidino, guanidino, hydrazino, amidazonyl, C₁₋₄alkylamino, C₁₋₄dialkylamino, halogen, perfluoro C₁₋₄alkyl, C₁₋₄alkyl, C₁₋₃alkoxy, nitro, carboxy, cyano, sulfuryl or hydroxyl), imidazoC14alkyl, substituted imidazol C14alkl (where the imidazole sustituents are independently selected from one or more of amino, amidino, guanidino, hydrazino, amidazonyl, C₁₋₄alkylamino, C₁₋₄dialkylamino, halogen, perfluoro C₁₋₄alkyl, C₁-

2.-4. (Canceled)

(Currently Amended) The compound of claim I, wherein A is -(C=O)-, B is -(CHR₄)-, D is -(C=O)-, E is -(ZR₆)-, G-is-XR₂-, and the compound has the following formula (IV):

$$\begin{array}{c} R_{1} \\ W \\ R_{7} \\ N_{7} \\ N_{$$

wherein R₁, R₂, R₄, R₆, R₇, W, X and Z are as defined in claim 1.

 (Currently Amended) <u>A compound having The compound of claim-5</u>, wherein the compound has the following formula (VI):

$$R_b$$
 NH O R_a (VI) O O X_1 X_2

wherein

R_n is a phenyl group; a substituted phenyl group having one or more substituents wherein the one or more substituents are independently selected from one or more of amino, amidino, guanidino, hydrazino, amidazonyl, C₁₋₄alkylamino, C₁₋₄dialkylamino, halogen, perfluoro C₁₋₄alkyl, C₁₋₃alkoxy, nitro, carboxy, cyano, sulfuryl, and hydroxyl groups; a benzyl group; a substituted benzyl group with one or more substituents where the one or more substituents where the one or more substituents are—independently selected from one or more of amino, amidino, guanidino, hydrazino, amidazonyl, C₁₋₄alkylamino, C₁₋₄dialkylamino, halogen, perfluoro C₁₋₄alkyl, C₁₋₃alkoxy, nitro, carboxy, cyano, sulfuryl, and hydroxyl groups; or a bicyclic aryl group having 8 to 11 ring members, which may have 1 to 3 heteroatoms selected from nitrogen, oxygen or sulfur:

R_b is a monocyclic aryl group having 5 to 7 ring members, which may have 1 to 2 heteroatoms selected from nitrogen, oxygen or sulfur, and aryl ring in the compound-may have one or more substituents selected from a group consisting of halide, hydroxy, cyano, lower alkyl, and lower alkoxy groups;

 $R_c \ is \ a \ saturated \ or \ unsaturated \ C_{1-6} alkyl, C_{1-6} alkoxy, \ \underline{or} \ perfluoro \ C_{1-6} alkyl \ group;$ and

 X_1 , X_2 , and X_3 may be the same or different and independently selected from hydrogen, hydroxyl, and halide.

7. (Currently Amended) The compound of claim 6, wherein

 R_a is a phenyl group; a substituted phenyl group having one or more substituents wherein the one or more substituents are-independently selected from one or more of amino, amidino, guanidino, hydrazino, amidazonyl, C_{1-4} alkylamino, C_{1-4} dialkylamino, halogen, perfluoro C_{1-4} alkyl, C_{1-4} alkyl, C_{1-4} alkyl, C_{1-5} alkoxy, nitro, carboxy, cyano, sulfuryl, and hydroxyl groups; a benzyl group; a substituted benzyl group with one or more substituents where the one or more substituents are—independently selected from one or more of amino, amidino, guanidino, hydrazino, amidazonyl, C_{1-4} alkylamino, C_{1-4} dialkylamino, halogen, perfluoro C_{1-4} alkyl

C₁₋₃alkoxy, nitro, carboxy, cyano, sulfuryl, and hydroxyl groups; a naphthyl group; a quinolinyl group; or an isoquinolinyl group; and

 R_b is phenyl, pyridyl or piperidyl, all of which may be substituted with one or more substituents selected from a group consisting of halide, hydroxy, cyano, lower alkyl, and lower alkoxy groups.

8. (Currently Amended) The compound of claim 6, wherein

 R_a is a phenyl group; a substituted phenyl group having one or more substituents wherein the one or more substituents are-independently selected from one or more of amino, amidino, guanidino, hydrazino, amidazonyl, C_{1-4} alkylamino, C_{1-4} dialkylamino, halogen, perfluoro C_{1-4} alkyl, C_{1-3} alkoxy, nitro, carboxy, cyano, sulfuryl, and hydroxyl groups; a benzyl group; a substituted benzyl group with one or more substituents where the one or more substituents are-independently selected from one or more of amino, amidino, guanidino, hydrazino, amidazonyl, C_{1-4} alkylamino, C_{1-4} dialkylamino, halogen, perfluoro C_{1-4} alkyl, C_{1-3} alkoxy, nitro, carboxy, cyano, sulfuryl, and hydroxyl groups; or a naphthyl group; and

R_b is phenyl, which may be substituted with one or more substituents selected from a group consisting of halide, hydroxy, cyano, lower alkyl, and lower alkoxy groups.

9.-11. (Canceled)

 (Currently Amended) A pharmaceutical composition comprising a compound according to claim 1 or to claim 6 and pharmaceutically acceptable carrier.

13.-42. (Canceled)